10/748,615

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	569	514/564.ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:32
L2	5	1 and arginine.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:32
L3	12	1 and lysine.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:33
L4	1435	carnitine.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:33
L5	328	4 and composition.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:35
L6	895	514/58.ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:35
L7	569	514/564.ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:35
L8	1062	514/546.ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:36
L9	1113	514/547.ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:36

L10	14	6 and (7 8 9)	US-PGPUB; USPAT;	OR	ON	2006/05/12 21:36
			USOCR; EPO; JPO; DERWENT			
S1	82	arginine same pyrrolidone same carboxylate	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 16:41
S3	3070	lysine.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 17:51
S4	0	514/ccls.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON .	2005/07/08 17:51
S5	5675	hydrochloride.ti.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 17:52
S6	0	S3 and S5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 17:52
S10	1	cortisol same maltodextrin	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 18:44
S11	58	cortisol and maltodextrin	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 18:45
S12	4	S11 and hgh	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/08 18:45
S13	1	arginine-2-pyrrolidone-5-carboxylate	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2005/07/21 15:29

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S14	22	arginine adj pyroglutamate	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:32
S15	0	arginine adj 5-oxo-proline	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:33
S16	10	"914342"	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:33
S17	0	wo9801474	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:33
S18	14	"9801474"	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:37
S19	2	"4388325".pn.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/09 14:37
S20	2	cortisol adj suppressant	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 15:55
S21	0	cortisol adj suppressent	US-PGPUB; USPAT; USOCR;	OR	ON	2006/05/12 15:55
			EPO; JPO; DERWENT			
S22	0	cprtisol same (amino adj acid)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 15:55
S23	527	cortisol same (amino adj acid)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 15:56

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S24	5	S23 same (hgh somatotropin)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:00
S25	59	S23 same uptake	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 15:58
S26	3	S23 same reuptake	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 15:58
S27	0	maltodextrin same (hgh somatotropin)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:00
S28	0	carnatine same (hgh somatotropin)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:00
S29	6	carnitine same (hgh somatotropin)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:02
S30	30	hgh same cortisol	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:09
S31	2	maltodextrin same cortisol	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:24
S32	69	hgh same anabolic	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:24
S33	49	cortisol same catabolic	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:34

			sco. y			
S34	0	S32 and S33	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:25
S35	8	S33 and maltodextrin	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:26
S36	4	hgh same (nutritional adj supplement)	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:44
S37	58	cortisol same carnitine	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:45
S38	0	acetyl adj carnatine	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:46
S39	220	acetyl adj carnitine	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:24
S40	82	S39 and muscle	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:46
S41	8	S39 same muscle	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:47
S42	5	S39 and cortisol	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:51
S43	9	S39 and catabolic	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 16:53

			<u>-</u>			
S44	73	maltodextrin and cortisol	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:12
S45	73	maltodextrin same supplement same nutritional	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:13
S46	6842	nutritional adj supplement	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:14
S47	4	S46 same hgh	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:15
S48	3	S46 same S39	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:16
S49	1	S46 same carnatine	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:24
S50	233	S46 same carnitine	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:25
S51	1534	mitochondria? same muscle	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 17:26
S52	96	mitochondria? near2 muscle	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/05/12 21:32

10/748,615

pyrrolidone-5-carboxylate

Caution: A net charge appears to be present

5-oxo-proline

Caution: A net charge appears to be present

$$H_2N$$
 H_2N
 OH
 OH

arginine-2-pyrrolidone-5-carboxylate



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PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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=> => file reg COST IN U.S. DOLLARS

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.21 0.21

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E5
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E6
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E11
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=> file caplus
COST IN U.S. DOLLARS
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FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 0.88 1.09 FILE 'CAPLUS' ENTERED AT 14:20:51 ON 09 MAY 2006
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FILE COVERS 1907 - 9 May 2006 VOL 144 ISS 20 FILE LAST UPDATED: 8 May 2006 (20060508/ED)
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=> d his

L2

L3

L4

(FILE 'HOME' ENTERED AT 14:19:38 ON 09 MAY 2006)

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FILE 'REGISTRY' ENTERED AT 14:19:44 ON 09 MAY 2006
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E 657-27-2/RN

L1 1 S E3

E 64855-91-0/RN

1 S E3

E 3040-38-8/RN

1 S E3

E 9050-36-6/RN

1 S E3

FILE 'CAPLUS' ENTERED AT 14:20:51 ON 09 MAY 2006

=> s 11 and 12

1118 L1

13 L2

L5 1 L1 AND L2

=> s 13 or 14

911 L3

3508 L4

4417 L3 OR L4

=> s 15 and 16

L7 1 L5 AND L6

=> d bib abs 17

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN

2005:570531 CAPLUS

DN 143:83512

TI Nutritional supplement for enhancing the production and effect of natural human growth hormone

IN Nerenberg, Arnold P.

PA USA

AN

SO U.S. Pat. Appl. Publ., 7 pp.

CODEN: USXXCO

DT Patent

LA English

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO. DATE

ΡÌ US 2005143343 Α1 20050630 US 2003-748615 20031230

PRAI US 2003-748615 20031230

A nutritional supplement for enhancing the production and effect of natural human growth hormone includes: L-arginine-2-pyrrolidone-5-carboxylate in an amount of about 500 mg to about 10 g; L-lysine-HCl in an amount of about 500 mg to about 10 g; and a cortisol suppressant including at least one of acetyl-L-carnitine in an amount of about 1 g to about 10 g and maltodextrin in an amount of about 1 g to about 10 q

=> s 1213 L2

=> d ibib abs 1-13 18

ANSWER 1 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:570531 CAPLUS

DOCUMENT NUMBER: 143:83512

TITLE: Nutritional supplement for enhancing the production

and effect of natural human growth hormone

INVENTOR(S): Nerenberg, Arnold P.

PATENT ASSIGNEE(S):

SOURCE: U.S. Pat. Appl. Publ., 7 pp.

CODEN: USXXCO

DOCUMENT TYPE: Patent English LANGUAGE:

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

AUTHOR(S):

PATENT NO. KIND DATE APPLICATION NO. DATE ____ _____ A1 20050630 US 2003-748615 US 2005143343 20031230 US 2003-748615 PRIORITY APPLN. INFO.: 20031230

A nutritional supplement for enhancing the production and effect of natural human growth hormone includes: L-arginine-2-pyrrolidone-5-carboxylate in an amount of about 500 mg to about 10 g; L-lysine-HCl in an amount of about 500 mg to about 10 g; and a cortisol suppressant including at least one of acetyl-L-carnitine in an amount of about 1 g to about 10 g and maltodextrin in an amount of about 1 g to about 10 g.

ANSWER 2 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2005:485667 CAPLUS

DOCUMENT NUMBER: 143:165983

Ligand-Based Virtual Screening and in Silico Design of TITLE:

New Antimalarial Compounds Using Nonstochastic and

Stochastic Total and Atom-Type Quadratic Maps

Marrero-Ponce, Yovani; Iyarreta-Veitia, Maite; Montero-Torres, Alina; Romero-Zaldivar, Carlos; Brandt, Carlos A.; Avila, Priscilla E.; Kirchgatter,

Karin; Machado, Yanetsy

Department of Pharmacy, Faculty of Chemical Pharmacy CORPORATE SOURCE:

and Department of Drug Design, Chemical Bioactive Center, Central University of Las Villas, Santa Clara,

Villa Clara, 54830, Cubá

Journal of Chemical Information and Modeling (2005), SOURCE:

45(4), 1082-1100

CODEN: JCISD8; ISSN: /1549-9596

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal LANGUAGE: English

CASREACT 143:165983 OTHER SOURCE(S):

Malaria has been one of the most significant public health problems for centuries. It affects many tropical and subtropical regions of the world. The increasing resistance of Plasmodium spp. to existing therapies has heightened alarms about malaria in the international health community. Nowadays, there is a pressing need for identifying and developing new drug-based antimalarial therapies. In an effort to overcome this problem, the main purpose of this study is to develop simple linear discriminant-based quant. structure-activity relation (QSAR) models for

the classification and prediction of antimalarial activity using some of the TOMOCOMD-CARDD (TOpol. Mol. COMputer Design-Computer Aided "Rational" Drug Design) fingerprints, to enable computational scréening from virtual combinatorial datasets. In this sense, a database of 1562 organic chems. having great structural variability, 597 of them antimalarial agents and 965 compds. having other clin. uses, was analyzed and presented as a helpful tool, not only for theor. chemists but also for other researchers in this area. This series of compds. was py6cessed by a k-means cluster anal. to design training and predicting sers. Afterward, two linear classification functions were derived to discriminate between antimalarial and nonantimalarial compds. The models (including nonstochastic and stochastic indexes) correctly classify more than 93% of the compound set, in both training and external prediction datasets. They showed high Matthews' correlation coeffs., 0.889 And 0.866 for the training set and 0.855 and 0.857 for the test one. The models' predictivity was also assessed and validated by the random removal of 10% of the compds. to form a new test set, for which predictions were made using the models. overall means of the correct classification for this process (leave group 10% full-out cross validation) using the equations with nonstochastic and stochastic atom-based quadratic fingerprints were 93.93% and 92.77%, resp. The quadratic maps-based TOMOCOMD-CARDD approach implemented in this work was successfully compared with four of the most useful models for antimalarials selection reported to date. The developed models were then used in a simulation of a virtual search for Ras FTase (FTase = farnesyltransferase) inhibitor's with antimalarial activity; 70% and 100% of the 10 inhibitors used in this virtual search were correctly classified, showing the ability of the models to identify new lead antimalarials. Finally, these two QSAR models were used in the identification of previously/unknown antimalarials. In this sense, three synthetic intermediaries of quinolinic compds. were evaluated as active/inactive ones using the developed models. The synthesis and biol. evaluation of these chems. against two malaria strains, using chloroquine as a reference, was performed. An accuracy of 100% with the theor. predictions was observed Compound 3 showed antimalarial activity, being the first report of an arylaminomethylenemalonate having such behavior. This result opens a door to a virtual study considering a higher variability of the structural core already evaluated, as well as of other chems. not included in this study. We conclude that the approach described here seems to be a promising QSAR tool for the mol. discovery of novel classes of antimalarial drugs, which may meet the dual challenges posed by

drug-resistant parasites and the rapid progression of malaria illnesses.

REFERENCE COUNT: 111 THERE ARE 111 CITED REFERENCES AVAILABLE FOR

THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE
FORMAT

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ANSWER 3 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER:
                            2005:371025 CAPLUS
DOCUMENT NUMBER:
                            142:417205
                            Nitric oxide topical technology
TITLE:
INVENTOR(S):
                            Leitman, Lorn; Barni, Gustavo
PATENT ASSIGNEE(S):
                            U.S. Pat / Appl. Publ., 5 pp.
SOURCE:
                            CODEN: VSXXCO
DOCUMENT TYPE:
                            Patent,
LANGUAGE:
                            English
FAMILY ACC. NUM. COUNT:
PATENT INFORMATION:
     PATENT NO.
                            KÁND
                                   DATE
                                                 APPLICATION NO.
                                                                           DATE
                             l _ _ _
     US 2005090545
                            A1
                                   20050428
                                                 US 2003-691025
                                                                           20031022
PRIORITY APPLN. INFO.:
                                                 US 2003-691025
     The invention enables application of L-arginine made with arginine
     \alpha-ketoglutarate, arginine pyroglutamate, arginine ketoisocaproate and ornithine \alpha-ketoglutarate to work together and sep. along with
     other ingredients, topica/lly. It is intended to improve the process by
     which these compds. work/in the human organism. The invention is based on
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the facts that L-arginine is: (1) the immediate precursor of the endogenous vasodilator found in the arterial blood vessels,

endothelium-derived relaxing factor (EPRF), required for protein synthesis and, depending on the organism's needs, can either be metabolized to support glucose synthesis or catabolized to produce energy. This simplified noninvasive application surpasses gastrointestinal digestion so the compds. will not degrade and favoring absorption into the circulatory system thus enhancing the compound effects and at much lower dosage levels than oral or parenteral administration. A composition comprises L-arginine to be used topically that creates increases in muscle size, strength, endurance and power output. Thus, a composition consisted essentially of L-arginine α -ketoglutarate dihydrate 1000, potassium bicarbonate 100, sodium bicarbonate 100, glycine 100, iso-Pr myristate 800, water 700, and vitamin skin smoother with corn 2000 mg.

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L8 ANSWER 4 OF 13 CAPLUS COPYRIGHT 2006, ACS on STN
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ACCESSION NUMBER: 2005:244333 CAPLÓS

DOCUMENT NUMBER: 143:307

AUTHOR (S):

TITLE: Atom, atom-type, and total nonstochastic and

stochastic quadratic fingerprints: a promising approach for modeling of antibacterial activity Marrero-Ponce, Yovani; Medina-Marrero, Ricardo;

Torrens, Francisco; Martinez, Yamile; Romero-Zaldivar,

Vicente; Castro, Eduardo A.

CORPORATE SOURCE: Department/of Pharmacy, Faculty of Chemical-Pharmacy,

Central University of Las Villas, Santa Clara, 54830,

Cuba

SOURCE: Bioorganic Medicinal Chemistry (2005), 13(8),

2881-2899

CODEN: BMECEP; \SSN: 0968-0896

PUBLISHER: Elsevier Ltd.

DOCUMENT TYPE: Journal LANGUAGE: English

The Topol. Mol. Computer Design (TOMOCOMD-CARDD) approach has been introduced for the classification and design of antimicrobial agents using computer-aided mol. design. For this propose, atom, atom-type, and total quadratic indexes have been generalized to codify chemical structure information. In this sense, stochastic quadratic indexes have been introduced for the description of the mol. structure. These stochastic fingerprints are based on a simple model for the intramol. movement of all valence-bond electrons. In this work, a complete data set containing 1006 antimicrobial agents is collected and presented. Two structure-based antibacterial activity classification models have been generated. The models (including nonstochastic and stochastic indexes) classify correctly more than 90% of 1525 compds. in training sets. These models permit the correct classification of 92.28% and 89.31% of 505 compds. in an external test sets. The approach, also, satisfactorily compares with respect to nine of the most useful models for antimicrobial selection reported to date. Finally, a virtual screening of 87 new compds. reported in the anti-infective field with antibacterial activities is developed showing the ability of the models to identify new leads as antibacterial.

REFERENCE COUNT: 91 THERÉ ARE 91 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 5 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:991310 CAPLUS

DOCUMENT NUMBER: 140:31162

TITLE: Use of an agent mimicking dopachrome tautomerase

(Trp-2) activity as protective agent for hair follicle

melanocytes and uses thereof

INVENTOR(S): Commo, Stephane; Gaillard, Olivier; Bernard, Bruno

PATENT ASSIGNEE(S): L'oreal, Fr. SOURCE: PCT Int. App

PCT Int. Appl., 32 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: French

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

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WO 2003103616
                                               WO 2003-FR1729
                                                                        20030610
                            A2
                                  20031218
     WO 2003103616
                            Α3
                                  20040415
             AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
              CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
              GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
              LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM,
              PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT,
              TZ, UA, UG, US, UZ, VC, VN, YÚ, ZA, ZM, ZW
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,
              KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES,
              FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
              BF, BJ, CF, CG, CI, CM, ÇÁ, GN, GQ, GW, ML, MR, NE, SN, TD, TG
     FR 2840531
                            A1
                                  20031/212
                                               FR 2002-7137
     FR 2840531
                            B1
                                  2004/1029
     CA 2487945
                                  20031218
                                               CA 2003-2487945
                            AΑ
                                                                        20030610
     AU 2003255653
                                  20031222
                                               AU 2003-255653
                            Α1
                                                                        20030610
     EP 1515688
                            A2
                                  20050323
                                               EP 2003-757134
                                                                        20030610
             AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
              IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK
     JP 2006512285
                            T2
                                  20060413
                                               JP 2004-510737
                                                                        20030610
     US 2005186233
                            A1
                                  20050825
                                               US 2004-9153
                                                                        20041213
PRIORITY APPLN. INFO.:
                                               FR 2002-7137
                                                                     A 20020611
                                               US 2002-389708P
                                                                     P
                                                                        20020619
                                                                     W 20030610
                                               WO 2003-FR1729
     The invention concerns the cosmeti/c use of an agent mimicking dopachrome
AB
     tautomerase (Trp-2) activity as ptotective agent for hair follicle
     melanocytes and its use, in particular for fighting against canities.
     invention also concerns specific/cosmetic compns. for fighting against
     canities comprising in a cosmetically acceptable medium at least an agent mimicking dopachrome tautomerase (Trp-2) activity and their uses. The
     invention further concerns a method for treating canities and a method for
     preserving natural pigmentation of gray or white hair and/or hairs by
     applying a cosmetic composition comprising at least one agent mimicking
     dopachrome tautomerase activity Finally, the invention concerns a method for identifying at least one agent mimicking dopachrome tautomerase
     (Trp-2) activity and a method for evaluating its cytoprotective activity.
     A hair lotion contained dopachrome tautomerase 0.5, propylene glycol 20,
     ethanol 30 and water q.s. 100 g.\
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L8 ANSWER 6 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2003:971575 CAPLUS

DOCUMENT NUMBER: 140:31172

TITLE: Cosmetic composition containing an agent mimicking the

activity of dopachrome tautomerase (Trp-2) to prevent

hair whitening

INVENTOR(S): Commo, Stephane; Gaillard, Olivier; Bernard, Bruno

PATENT ASSIGNEE(S): L'oreal, Fr.

SOURCE: Fr. Demande, 39 pp.

CODEN: FRXXBL

DOCUMENT TYPE: Patent LANGUAGE: French

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT	NO.			KIN	D :	DATE			APPL	ICAT	ION 1	NO.		Dž	ATE	
					-									-		
FR 2840	531			A 1		2003	1212		FR 2	002-	7137			20	0020	511
FR 2840	531			В1		2004	1029									
CA 2487	945			AA		2003	1218		CA 2	003-	2487	945		20	0030	510
WO 2003	1036	16		A2		2003	1218	1	WO 2	003-	FR17	29		20	0030	510
WO 2003	1036	16		A3		2004	0415									
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						MD,										-
	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	ТJ,	TM,	TN,	TR,	TT,
	TZ,	UA,	UG,	US,	UZ,	VC,	VN,	YU,	ZA,	ZM,	ZW	•	•	•	•	•
RW:	GH,	•		•	•	•	•	•	•	•		ZM.	ZW.	AM.	AZ.	BY.
	•		-			TM,	•	•		•	•	•	•		•	•

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FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
               BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
      AU 2003255653

      20031222
      AU 2003-255653
      20030610

      20050323
      EP 2003-757134
      20030610

                              A1
                              A2
          R: AT, BE, CH, DE, DK, ES, FR, GB, GR, LT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK
      JP 2006512285 T2 20060413
US 2005186233 A1 20050825
                                                 JP ∕2004-510737
                                                    US 2004-9153
FR 2002-7137
PRIORITY APPLN. INFO.:
                                                                         A 20020611
                                                    US 2002-389708P P 20020619
WO 2003-FR1729 W 20030610
      A cosmetic composition to fight against the/hair whiteness contains an agent
      mimicking the activity of dopachrome tautomerase (Trp-2). The invention
      refers moreover to a method for identifying an agent mimicking the
      activity of Trp-2. Expression of Trp-2 in melanocytes from human hair
      follicles and epidermis is studied. A hair lotion contained Trp-2 0.5,
      propylene glycol 20, ethanol 30, and water q.s. 100 g.
                             22
                                    THERE ARE 22 CITED REFERENCES AVAILABLE FOR THIS
REFERENCE COUNT:
                                    RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT
     ANSWER 7 OF 13 CAPLUS COPYRIGHT/2006 ACS on STN
ACCESSION NUMBER: 2000:116751 /CAPLUS
DOCUMENT NUMBER:
                             132:156552
TITLE:
                           Deodorant and/or moisturizing cosmetic composition
                          containing an orthophosphoric acid ester and polyacrylate
Mucci, Paolo; Meucci, Sandro; Ceccarelli, Luigi
Societa Italo-Britannica L. Manetti-H. Roberts & C.
INVENTOR(S):
PATENT ASSIGNEE(S):
                             S.p.A., Italy
                             Eur. Pat. Appl., 10 pp.
SOURCE:
                             CODEN: EPXXDW
DOCUMENT TYPE:
                             Patent
LANGUAGE:
                             English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:
                                                  APPLICATION NO.
                           KIND DATE
     PATENT NO.
                                     20000216 EP 1999-830439
     EP 979644
                             ----
                                                   -----
                              A1
                                                                               19990708
          R: AT, BE, CH, DE, DK, ES, \FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
               IE, SI, LT, LV, FI, RO
                              B1 20000720 IT 1998-MI1895 19980813
IT 1998-MI1895 A 19980813
      IT 1302023
PRIORITY APPLN. INFO.:
     A deodorant and/or moisturizing cosmetic composition formed by an emulsion oil
     in water, comprising: (a) a combination of two or more deodorant and/or moisturizing active ingredients, (b) an orthophosphoric acid ester in an amount comprised between 0.3 and 7%, and (c) a water soluble resin consisting
     of a homopolymer of the acrylic acid in an amount comprised between 0.05%
      and 1.5%, said composition having a viscosity at 20°C comprised between
      800 and 10000 caps. The main advantage of the cosmetic composition according
      to the invention consists in that it is a cream and, at the same time, is
     vaporizable, as the viscosity is always comprised in the above stated range. A vaporizable cream contained water 74.876, vaseline oil 8.000, propylene glycol 5.000, Myritol 312 5.000, Hostaphat KL 340 N 3.000,
     Cosmacol ELI 1.000, perfume 1.000, farnesol 0.6000, vitamin E acetate
      0.500, Germaben II 0.500, Carborol 5/984 0.300, sodium hydroxide 0.124,
     disodium EDTA, and BHT 0.050%.
REFERENCE COUNT:
                                    THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS
                             4
                                    RECORD! ALL CITATIONS AVAILABLE IN THE RE FORMAT
     ANSWER 8 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN
                             1998:65929 CAPLUS
ACCESSION NUMBER:
DOCUMENT NUMBER:
                             128:128292
TITLE:
                             Preparation of somatostatin derivatives and their
                             combinations with amino acids or oligopeptides for
                             promoting body growth
INVENTOR (S):
                             Volpato, Ivo; Bizzini, Bernard; Grabitz, Ernst
                             Bernhard
PATENT ASSIGNEE(S):
                             Dox-Al Italia S.P.A., Italy
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SOURCE:
                          PCT Int. Appl., 39 pp.
                          CODEN: PIXXD2
DOCUMENT TYPE:
                          Patent
LANGUAGE:
                          English
FAMILY ACC. NUM. COUNT:
PATENT INFORMATION:
                         KIND DATE
                                               APPLICATION NO.
     PATENT NO.
     -----
                          ----
                                              -----<del>\</del>-------
     WO 9801474
                          A2 19980115
                                               WO 1997-EP3605
     WO 9801474
                           A3
                                  19980409
         W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ,
             LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL,
             PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ,/TM, TR, TT, UA, UG, US,
             UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD,/RU, TJ, TM
         RW: GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE/ CH, DE, DK, ES, FI, FR,
             GB, GR, IE, IT, LU, MC, NL, PT, SE, B_{\overline{F}}^{\prime}, BJ, CF, CG, CI, CM, GA,
             GN, ML, MR, NE, SN, TD, TG
     AU 9736930
                                19980202 AU 1997-36930
19990512 EP 1997-933655
                                              AU 1/997-36930
                           A1
                                19980202
     EP 914342
                           A2
            AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
             IE, FI
PRIORITY APPLN. INFO.:
                                               IT 1996-MI1408
                                               WO 1997-EP3605
                          MARPAT 128 128292
OTHER SOURCE(S):
     ANSWER 9 OF 13 CAPLUS COPYRIGHT 2006 AGS on STN
ACCESSION NUMBER:
                          1988:124398 CAPLUS/
DOCUMENT NUMBER:
                          108:124398
                          Effect of pyroglutamic acid stereoisomers on ECS and
TITLE:
```

Chemical modified somatostatins administered to man and animals can significantly increase the body growth rate and index. Furthermore, L-amino acids (in particular arginine pyroglutamate) or oligopeptides, administered orally in combination with the claimed derivs. of chemical modified somatostatins, produce a\synergic effect on body growth. Thus, polymerized somatostatin was prepared and shown to significantly increase animal growth (178.5 vs. 144.6% for the placebo after 21 days).

scopolamine-induced memory disruption and brain

19970708

19970708

19970708

19960708

19970708

Α

W

acetylcholine levels in the rat

AUTHOR(S): Spignoli, G.; Magnani, M.; Giovannini, M. G.; Pepeu,

CORPORATE SOURCE: Dep. Preclin./Clin. Pharmacol., Univ. Florence,

Florence, 501/34, Italy

SOURCE: Pharmacological Research Communications (1987),

19(12), 901-12

CODEN: PLRCAT, ISSN: 0031-6989

DOCUMENT TYPE: Journal LANGUAGE:

English In rats, the acquisition of a passive avoidance conditioned response was disrupted by electroconvulsive shock (ECS) or scopolamine administration. DL-Pyroglutamic acid (DL-PCA) 500 and 1000 mg/kg prevented both the ECS and scopolamine-induced amnesia. Afginine alone was ineffective. Scopolamine brought about a 52 and/39% decrease, resp., in cortical and hippocampal acetylcholine (ACh) levels. DL-PCA 500 and 1000 mg/kg also prevented the decrease in brain ACh level. When the 2 isomers were studied sep., D-PCA was more effective than L-PCA and antagonized scopolamine-induced amnesia at/250 and 500 mg/kg. DL-PCA appears to be active on cortical and hippocampal cholinergic mechanisms and, like other 2-oxopyrrolidone derivs., has cognition-enhancing properties.

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ANSWER 10 OF 13 CAPLUS COPYRIGHT 2006 AOS on STN
ACCESSION NUMBER:
                        1988:106414 CAPLUS
DOCUMENT NUMBER:
                        108:106414
```

Effects of pyroglut/amic acid on learning and memory TITLE:

processes of the rat

AUTHOR (S): Drago, F.; Continella, G.; Valerio, C.; D'Agata, V.;

Astuto, C.; Spadafo, F.; Scapagnini, U.

CORPORATE SOURCE: Med. Sch., Univ. Catania, Catania, 95125, Italy SOURCE: Acta Therapeutica (1987), 13(6), 587-94

CODEN: ACTTDZ; ISSN: 0378-0619

DOCUMENT TYPE: Journal LANGUAGE: English

AB The arginine salt of pyroglutamic acid (PCA) was administered subchronically to male rats (i.p. injection of 0.1 and 1 g/kg/day for 15 days). The PCA did not modify the rate of acquisition of the pole-jumping response, but inhibited its extinction. The dose of 1 g/kg was more potent than 0.1 g/kg in this respect. In addition, in the passive avoidance task, treatment with PCA was followed by an improvement of avoidance retention. Both 24 and 48 h after the learning trial, PCA-treated rats showed better memory retention than control animals.

L8 ANSWER 11 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1982:515326 CAPLUS

DOCUMENT NUMBER: 97:115326

TITLE: L-Arginine-DL-pyroglutamate as an agent with an effect

on a neuroendocrine area

INVENTOR(S): Orzalesi, Giovanni

PATENT ASSIGNEE(S): Societa Italo-Britannica L. Manetti-H. Roberts e C.,

Italy

SOURCE: Ger. Offen., 20 pp.

CODEN: GWXXBX

DOCUMENT TYPE: Patent LANGUAGE: German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	AP	PLICATION NO.		DATE
DE 3125512	A1	19820506	DE	1981-3125512		19810629
US 4388325	A	19830614	-	1981-275774		19810622
ZA 8104255	A	19820728	ZA	1981-4255		19810624
JP 57098213	A2	19820618	JP	1981-99793		19810629
CA 1168985	A1	19840612	CA	1981-380896		19810630
PRIORITY APPLN. INFO.:			IT	1980-49111	Α	19800630
AB L-arginine DL-pyro	glutamat	e [64855-91	L-0]	Can increase t	he	

L-arginine DL-pyroglutamate [64855-91-0] Can increase the sexual activity of elderly male mammals when administered daily at 200-4500 mg orally or 200-1200 mg parenterally. The compound increases dopaminergic tone. Expts. on the response of learning and sexual behavior of young and old rats and on the sexual behavior of men under 40 and over 60 yr old to the peptide are described. There was no effect on the sexual activity of young men, but that of older men was significantly increased. Capsules were prepared containing peptide 1000, modified starch 98, and Mg stearate 2 mg.

ANSWER 12 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1978:517679 CAPLUS

DOCUMENT NUMBER: 89:117679

TITLE: Chemicopharmaceutic and pharmacodynamic

characteristics of pyrglutargine

AUTHOR(S): Selleri, R.; Orzalesi, G.; Innocenti, F.; Volpato, I.;

Bisagno, T.

CORPORATE SOURCE: Lab. Ric. Farmacobiol., Soc. Italo-Britannica,

Calenzano, Italy

SOURCE: Bollettino Chimico Farmaceutico (1977), 116(12),

735-43

CODEN: BCFAAI; ISSN: 0006-6648

DOCUMENT TYPE: Journal LANGUAGE: Italian

LANC

$$0 \xrightarrow{\text{H}} \begin{array}{c} \text{NH} & \text{NH}_2 \\ \text{II} & \text{I} \\ \text{II} & \text{I} \\ \text{CO}_2 - & \text{+H}_3 \text{NCNH} (\text{CH}_2) \text{ 3CHCO}_2 \text{H} \\ \end{array}$$

Reaction of L-arginine with DL-pyroglutamic acid for a few min in boiling water, followed by evaporation of the solvent and crystallization of the product from 95% EtOH-H2O (7:3), yielded pyrglutargine (I) [64855-91-0], m. 216-17°, $[\alpha]$ 250 +11.9° (c 7.00, H2O), conductivity 54.4 S + mol-1. I is probably a pair of LL- and LD-diastereoisomers. IR, NMR, and mass spectra of I and of the 2 reactant amino acids are also illustrated and discussed. A review is given of previously published results on the pharmacol. of the compound which showed it to affect the central nervous system, inhibiting the effects of neurodepressant drugs and facilitating the learning of some specialized behavior. The formulation of capsules and liquid solns. of I is also described.

ANSWER 13 OF 13 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1978:15861 CAPLUS

DOCUMENT NUMBER: 88:15861

TITLE: Chemistry and pharmacology of arginine pyroglutamate.

Analysis of its effects on the CNS

AUTHOR (S): Provenzano, P. M.; Brucato, A.; Gianguzza, S.;

Coppola, A.; Orzalesi, G.; Selleri, R.; Innocenti, F.;

Volpato, I.

Dep. Toxicol., Univ. Palermo, Palermo, Italy 🧨 CORPORATE SOURCE:

Arzneimittel-Forschung (1977), 27(8), 1553-7 SOURCE:

CODEN: ARZNAD; ISSN: 0004-4172

DOCUMENT TYPE: Journal English

LANGUAGE:

CO2H H2NCNH (CH2) 3

AB The effects of arginine pyroglutamate (I) [64855-91-0] on the central nervous system in rats were studied. I antagonized the general anesthesia produced by Na pentobarbital [57-33-0] even in the presence of medazepam [2898-12-6] and flurazepam [17617-23-1]. I also attenuated the decrease in locomotor behavior induced by pentobarbital and the benzodiazepines. I did not alter the sound discrimination capacity at fixed intervals nor did it influence the learning of a sound discrimination at varied intervals. Learning was moderately accelerated by I in temporal discrimination and conditioned avoidance response tests. Neither arginine nor pyroglutamate had any effect when given alone. I appears to block the effect of central nervous system depressants without affecting normal behavior.

148,615

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=> e l-arginine-2-pyrrolidone-5-carboxylate/cn
              1
                    L-ARGININE-2-D/CN
E2
              1
                    L-ARGININE-2-D, MONOHYDROCHLORIDE/CN
E3
              0 --> L-ARGININE-2-PYRROLIDONE-5-CARBOXYLATE/CN
E4
              1
                    L-ARGININE-3,4,5,5-T4/CN
E5
              1
                    L-ARGININE-3,4-T2/CN
E6
              1
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E7
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                    L-ARGININE-4,4-D2/CN
E8
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                    L-ARGININE-4,5-T2/CN
E9
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                    L-ARGININE-4-D, ERYTHRO-/CN
E10
              1
                    L-ARGININE-4-D, THREO-/CN
E11
              1
                    L-ARGININE-4-NITROANILIDE HYDROCHLORIDE/CN
E12
                    L-ARGININE-5,5-D2, 3-HYDROXY-, ERYTHRO-/CN
=> e l-lysine hydrochloride/cn
                    L-LYSINE EXPORTER (SHEWANELLA ONEIDENSIS STRAIN MR-1 GENE SO
                    2865)/CN
F.2
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                    OLYMER/CN
E5
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                    R/CN
E6
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E7
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            1 L-LYSINE HYDROFLUORIDE/CN
1 L-LYSINE HYDROXAMATE HYDROCHLORIDE/CN
1 L-LYSINE HYDROXAMIC ACID/CN
E8
E9
E10
E11
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                    L-LYSINE ISOPROPYL ESTER/CN
E12
                  L-LYSINE L-GLUTAMATE/CN
=> s e3
              1 "L-LYSINE HYDROCHLORIDE"/CN
=> d 11
L1
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN
     657-27-2 REGISTRY
ED
     Entered STN: 16 Nov 1984
     L-Lysine, monohydrochloride (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Lysine, monohydrochloride, L- (8CI)
OTHER NAMES:
CN
    Darvyl
CN
     L-Gen
CN
     L-Lysine hydrochloride
CN
     Lyamine
CN
     Lysine hydrochloride
CN
     Lysine monohydrochloride
CN
     Lysion
     NSC 9253
CN
FS
     STEREOSEARCH
     305-76-0, 93394-22-0
DR
MF
     C6 H14 N2 O2 . Cl H
CI
     COM
LC
     STN Files:
                   AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DETHERM*, DIOGENES, EMBASE, GMELIN*, HSDB*,
       IFICDB, IFIPAT, IFIUDB, IPA, MRCK*, MSDS-OHS, PATDPASPC, PROMT, RTECS*,
       SPECINFO, TOXCENTER, USAN, USPAT2, USPATFULL
          (*File contains numerically searchable property data)
     Other Sources:
                       DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

$$NH_2$$
 HO_2C
 S
 $(CH_2)_4$
 NH_2

● HCl

CN

L-O-Acetylcarnitine

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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1067 REFERENCES IN FILE CA (1907 TO DATE)
29 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1067 REFERENCES IN FILE CAPLUS (1907 TO DATE)
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=> e acetyl-l-carnitine/cn
E1
             1
                   ACETYL-L-ALANYLGLYCYLGLYCINE METHYL ESTER/CN
E2
             1
                   ACETYL-L-ASPARTIC ACID/CN
E3
             1 --> ACETYL-L-CARNITINE/CN
             1
E4
                 ACETYL-L-CARNITINE ACID PHOSPHATE/CN
E5
             1
                   ACETYL-L-CARNITINE ACID SULFATE/CN
E6
             1
                  ACETYL-L-CARNITINE GLUCOSE PHOSPHATE/CN
E7
             1
                 ACETYL-L-CARNITINE GLYCEROPHOSPHATE/CN
                 ACETYL-L-CARNITINE LACTATE/CN
ACETYL-L-CARNITINE MAGNESIUM CITRATE/CN
E8
            1
E9
             1
E10
             1
                 ACETYL-L-CARNITINE METHANESULFONATE/CN
E11
                  ACETYL-L-CARNITINE OROTATE/CN
E12
                   ACETYL-L-CARNITINE TRICHLOROACETATE/CN
=> s e3
             1 ACETYL-L-CARNITINE/CN
=> d 12
L2
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN
     3040-38-8 REGISTRY
ED
     Entered STN: 16 Nov 1984
     1-Propanaminium, 2-(acetyloxy)-3-carboxy-N,N,N-trimethyl-, inner salt,
     (2R) - (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
     1-Propanaminium, 2-(acetyloxy)-3-carboxy-N,N,N-trimethyl-, inner salt,
CN
     Ammonium, (3-carboxy-2-hydroxypropyl)trimethyl-, hydroxide, inner salt,
     acetate, L- (8CI)
OTHER NAMES:
     (-)-Acetylcarnitine
CN
     (R)-Acetylcarnitine
CN
     Acetyl-L-(-)-carnitine
CN
     Acetyl-L-carnitine
CN
     Acetylcarnitine
CN
     ALCAR
CN
     L-Acetylcarnitine
CN
     L-Carnitine acetyl ester
```

```
CN
    Levocarnitine acetyl
CN
    Nicetile
CN
    O-Acetyl-L-carnitine
CN
    O-Acetylcarnitine
FS
    STEREOSEARCH
DR
    461-77-8, 541-68-4, 3624-25-7, 74832-89-6
MF
    C9 H17 N O4
CI
    COM
LC
    STN Files:
                 ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*,
      BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CEN,
      CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DRUGU, EMBASE, IMSCOSEARCH,
       IMSDRUGNEWS, IMSRESEARCH, IPA, MRCK*, PROMT, PROUSDDR, RTECS*,
       TOXCENTER, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
    Other Sources:
                      WHO
```

OTHER NAMES:

CN

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

857 REFERENCES IN FILE CA (1907 TO DATE)
19 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
858 REFERENCES IN FILE CAPLUS (1907 TO DATE)
3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

```
=> e l-arginine/cn
                   L-ARGININAMIDE-N-15N, L-VALYL-L-GLUTAMINYL-N-15N-L-ALANYL-L-
E1
             1
                   ASPARAGINYL-N-15N-L-ISOLEUCYL-L-ALANYLGLYCYL-L-HISTIDYLGLYCY
                   L-L-GLUTAMINYL-L-GLUTAMINYL-N-15N-L-VALYL-L-LEUCYL-L-ISOLEUC
                   YL-/CN
E2
             1
                   L-ARGININAMIDE-N2-15N, N-FORMYL-L-ALANYL-/CN
E3
             1 --> L-ARGININE/CN
                   L-ARGININE A-PHENOXYBUTYRATE/CN
E4
             1
E5
             1
                   L-ARGININE B-NAPHTHYLAMIDE/CN
E6
             1
                   L-ARGININE 2-NAPHTHYLAMIDE/CN
E7
             1
                   L-ARGININE 4'-ETHOXYAZOBENZENE-4-SULFONATE/CN
E8
             1
                   L-ARGININE 4-METHYLCOUMARYL-7-AMIDE/CN
E9
             1
                   L-ARGININE ACETYLSALICYLATE/CN
E10
                   L-ARGININE ACETYLSALICYLATE SALT (1:1)/CN
             1
E11
             1
                  L-ARGININE ACETYLSALICYLIC ACID SALT/CN
E12
                  L-ARGININE AMIDE/CN
=> s e3
L3
             1 L-ARGININE/CN
=> d 13
L3
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN
     74-79-3 REGISTRY
ED
     Entered STN: 16 Nov 1984
     L-Arginine (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
    Arginine, L- (8CI)
```

(S) -2-Amino-5-[(aminoiminomethyl)amino]pentanoic acid

```
CN
     Arginine
CN
     L-(+)-Arginine
CN
     L-\alpha-Amino-\delta-guanidinovaleric acid
CN
     L-Arg
CN
     L-Norvaline, 5-[(aminoiminomethyl)amino]-
CN
     L-Ornithine, N5-(aminoiminomethyl)-
CN
     NSC 206269
CN
     Pentanoic acid, 2-amino-5-[(aminoiminomethyl)amino]-, (S)-
FS
     STEREOSEARCH
     667422-95-9, 7004-12-8, 142-49-4
DR
MF
     C6 H14 N4 O2
CI
     COM
LC
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB,
       CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU,
       DETHERM*, DIOGENES, DRUGU, EMBASE, GMELIN*, HODOC*, HSDB*, IFICDB,
       IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, NIOSHTIC,
       PATDPASPC, PHAR, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE,
       TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, VETU
         (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

41313 REFERENCES IN FILE CA (1907 TO DATE)
1202 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
41367 REFERENCES IN FILE CAPLUS (1907 TO DATE)
6 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> e maltode	xtrin/	cn
E1	1	MALTOBIOURONOSIDE, BENZYL, METHYL ESTER, HEXAACETATE/CN
E2	2	MALTODECAOSE/CN
E3	1>	MALTODEXTRIN/CN
E4	1	MALTODEXTRIN 19/CN
E5	1	MALTODEXTRIN 24DE/CN
E6	1	MALTODEXTRIN ABC TRANSPORTER ATP-BINDING PROTEIN (MYCOPLASMA
		MOBILE STRAIN 163K GENE MALK)/CN
E7	1	MALTODEXTRIN ABC TRANSPORTER ATP-BINDING PROTEIN MALK (MYCOP
		LASMA PULMONIS STRAIN UAB CTIP GENE MYPU-6410)/CN
E8	1	MALTODEXTRIN ABC TRANSPORTER PERMEASE PROTEIN (MYCOPLASMA MO
		BILE STRAIN 163K GENE MALC)/CN
E9	1	MALTODEXTRIN ABC TRANSPORTER PERMEASE PROTEIN MALC (MYCOPLAS
		MA PULMONIS STRAIN UAB CTIP GENE MYPU-6390)/CN
E10	1	MALTODEXTRIN ABC TRANSPORTER PERMEASE PROTEIN MALD (MYCOPLAS
		MA PULMONIS STRAIN UAB CTIP GENE MYPU-6400)/CN
E11	1	MALTODEXTRIN ABC TRANSPORTER, PERMEASE PROTEIN (STREPTOCOCCU
		S PNEUMONIAE STRAIN TIGR4 GENE SP2109)/CN
E12	1	MALTODEXTRIN ABC TRANSPORTER, PERMEASE PROTEIN (STREPTOCOCCU
		S PNEUMONIAE STRAIN TIGR4 GENE SP2110)/CN

```
=> s e3
L4
             1 MALTODEXTRIN/CN
=> d 14
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN
     9050-36-6 REGISTRY
ED
     Entered STN: 16 Nov 1984
CN
     Maltodextrin (9CI) (CA INDEX NAME)
OTHER NAMES:
     Amidex DE 10
CN
     C Pharm 01980
CN
     C*De Light 01970
     C*deLight F 01970
CN
     C*deLight MD 01970
CN
CN
     C-Pur 01910
     C-PUR 01915
CN
     C-Pur 01921
CN
CN
     C-Sperse MD 01314
     Cerestar C*PUR 01915
CN
CN
     Cerestar PUR 01915
CN
     DE 2
CN
     Dextrin, malto
CN
     Dry Sweet
CN
     Fibersol 2(E)
CN
     Foodtex
CN
     Frodex 10
CN
     Frodex 20
CN
     Glucidex 12
CN
     Glucidex 17
     Glucidex 19
CN
     Glucidex 19FD
ĊN
CN
     Glucidex 2
CN
     Glucidex 21
CN
     Glucidex 2B
     Glucidex 39
CN
CN
     Glucidex 6
CN
     Glucidex IT 12
CN
     Glucidex IT 19
CN
     Glucidex IT 6
CN
     Instant N-Oil II
CN
     Instant Oil II
CN
     Instant Stellar
CN
     K 8
CN
     Lodex 10
CN
     Lodex 5
CN
     Lycadex 100
CN
     Lycadex 200
CN
     Lycatab
CN
     M 01960
CN
     M 040
CN
     Maldex 15
CN
     Maldex 150
     Maldex 20
CN
     Maldex 30
CN
CN
     Malta-Gran 10
CN
     Malta-Gran TG
CN
     Maltiva
     Maltodextrin 19
CN
CN
    Maltodextrin 24DE
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
     DISPLAY
DR
     126776-44-1, 126776-45-2, 127120-90-5, 54077-26-8, 104859-39-4,
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104859-43-0, 104859-45-2, 104859-47-4, 104859-49-6, 104859-62-3, 104859-75-8, 61008-41-1, 142583-82-2, 89750-26-5, 87090-11-7, 39283-25-5, 52769-80-9, 216252-89-0, 220857-34-1, 287179-53-7

MF Unspecified

CI PMS, COM, MAN

PCT Manual registration

LC STN Files: AGRICOLA, ANABSTR, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, PIRA, PROMT, TOXCENTER, USPAT2, USPATFULL

(*File contains numerically searchable property data)

Other Sources: DSL**, EINECS**, TSCA**

(**Enter CHEMLIST File for up-to-date regulatory information)

- *** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
- **PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**

3125 REFERENCES IN FILE CA (1907 TO DATE)

139 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

3125 REFERENCES IN FILE CAPLUS (1907 TO DATE)

(FILE 'HOME' ENTERED AT 16:47:13 ON 08 JUL 2005)

FILE 'REGISTRY' ENTERED AT 16:47:30 ON 08 JUL 2005 E L-ARGININE-2-PYRROLIDONE-5-CARBOXYLATE/CN E L-LYSINE HYDROCHLORIDE/CN L1 1 S E3 E ACETYL-L-CARNITINE/CN L2 1 S E3 E L-ARGININE/CN L3 1 S E3 E MALTODEXTRIN/CN

L4 . 1 S E3

10/748,615

Welcome to STN International! Enter x:x

LOGINID:ssspta1600txm

PASSWORD:

NEWS HOURS

NEWS LOGIN

NEWS IPC8

TERMINAL (ENTER 1, 2, 3, OR ?):2

Welcome to STN International NEWS Web Page URLs for STN Seminar Schedule - N. America NEWS 2 "Ask CAS" for self-help around the clock NEWS 3 DEC 23 New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/ USPAT2 NEWS 4 JAN 13 IPC 8 searching in IFIPAT, IFIUDB, and IFICDB NEWS 5 JAN 13 New IPC 8 SEARCH, DISPLAY, and SELECT enhancements added to TNPADOC NEWS 6 JAN 17 Pre-1988 INPI data added to MARPAT 7 JAN 17 IPC 8 in the WPI family of databases including WPIFV NEWS NEWS 8 JAN 30 Saved answer limit increased NEWS 9 FEB 21 STN AnaVist, Version 1.1, lets you share your STN AnaVist visualization results NEWS 10 FEB 22 The IPC thesaurus added to additional patent databases on STN NEWS 11 FEB 22 Updates in EPFULL; IPC 8 enhancements added NEWS 12 FEB 27 New STN AnaVist pricing effective March 1, 2006 NEWS 13 FEB 28 MEDLINE/LMEDLINE reload improves functionality NEWS 14 FEB 28 TOXCENTER reloaded with enhancements NEWS 15 FEB 28 REGISTRY/ZREGISTRY enhanced with more experimental spectral property data NEWS 16 MAR 01 INSPEC reloaded and enhanced NEWS 17 MAR 03 Updates in PATDPA; addition of IPC 8 data without attributes NEWS 18 MAR 08 X.25 communication option no longer available after June 2006 NEWS 19 MAR 22 EMBASE is now updated on a daily basis NEWS 20 APR 03 New IPC 8 fields and IPC thesaurus added to PATDPAFULL NEWS 21 APR 03 Bibliographic data updates resume; new IPC 8 fields and IPC thesaurus added in PCTFULL NEWS 22 APR 04 STN AnaVist \$500 visualization usage credit offered NEWS 23 APR 12 LINSPEC, learning database for INSPEC, reloaded and enhanced NEWS 24 Improved structure highlighting in FQHIT and QHIT display APR 12 in MARPAT NEWS 25 APR 12 Derwent World Patents Index to be reloaded and enhanced during second quarter; strategies may be affected NEWS EXPRESS FEBRUARY 15 CURRENT VERSION FOR WINDOWS IS V8.01a, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 19 DECEMBER 2005. V8.0 AND V8.01 USERS CAN OBTAIN THE UPGRADE TO V8.01a AT http://download.cas.org/express/v8.0-Discover/

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Take survey: http://www.zoomerang.com/survey.zgi?p=WEB2259HNKWTUW

Thank you in advance for your participation.

FILE 'HOME' ENTERED AT 12:18:30 ON 09 MAY 2006

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 12:18:51 ON 09 MAY 2006 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 9 May 2006 VOL 144 ISS 20 FILE LAST UPDATED: 8 May 2006 (20060508/ED)

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http://www.cas.org/infopolicy.html

=> s us 20050143343.pn 98845 US 3198926 USES 3290620 US (US OR USES) 0 20050143343 33567 PN 1998 PNS 35426 PN

> (PN OR PNS) 0 US 20050143343.PN

(US(W)20050143343(W)PN)

=> sel rn

I.1

=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 8.34 8.55

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 12:19:33 ON 09 MAY 2006
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STRUCTURE FILE UPDATES: 8 MAY 2006 HIGHEST RN 883439-06-3
DICTIONARY FILE UPDATES: 8 MAY 2006 HIGHEST RN 883439-06-3

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TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

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REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> s e1-e10

L3

1 107-35-7/BI (107-35-7/RN)1 12629-01-5/BI (12629-01-5/RN) 1 3040-38-8/BI (3040-38-8/RN) 1 50-23-7/BI (50-23-7/RN)1 502-65-8/BI (502-65-8/RN) 1 56-85-9/BI (56-85-9/RN)1 60-18-4/BI (60-18-4/RN)1 64855-91-0/BI (64855-91-0/RN) 1 657-27-2/BI (657-27-2/RN) 1 9050-36-6/BI (9050-36-6/RN)

(9050-36-6/RN)

10 (107-35-7/BI OR 12629-01-5/BI OR 3040-38-8/BI OR 50-23-7/BI OR 502-65-8/BI OR 56-85-9/BI OR 60-18-4/BI OR 64855-91-0/BI OR 657-27-2/BI OR 9050-36-6/BI)

```
ANSWER 1 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN
L3
     64855-91-0 REGISTRY
RN_
     Entered STN: 16 Nov 1984
     Proline, 5-oxo-, compd. with L-arginine (1:1) (9CI)
                                                            (CA INDEX NAME)
OTHER CA INDEX NAMES:
     DL-Proline, 5-oxo-, compd. with L-arginine (1:1)
     L-Arginine, compd. with 5-oxo-DL-proline (1:1)
     L-Arginine, compd. with 5-oxoproline (1:1) (9CI)
OTHER NAMES:
CN
     Arginine pyroglutamate
CN
     Pyrglutargine
FS
     STEREOSEARCH
MF
     C6 H14 N4 O2 . C5 H7 N O3
LC
     STN Files: BEILSTEIN*, BIOSIS, CA, CAPLUS, CHEMLIST, EMBASE, PHAR,
       PROMT, PS, TOXCENTER, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
                     EINECS * *
         (**Enter CHEMLIST File for up-to-date regulatory information)
     CM
          1
     CRN
          149-87-1
     CMF C5 H7 N O3
     CM
          2
     CRN
          74-79-3
     CMF C6 H14 N4 O2
Absolute stereochemistry.
     NH
            (CH<sub>2</sub>)<sub>3</sub>
                   NH_2
              13 REFERENCES IN FILE CA (1907 TO DATE)
               2 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
              13 REFERENCES IN FILE CAPLUS (1907 TO DATE)
=> d 2-10 13
L3
     ANSMER 2 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN
     12629-01-5 REGISTRY
RN
     Entered STN: 16 Nov 1984
ED
     Somatotkopin (human) (9CI)
CN
                                 (CA INDEX NAME)
OTHER NAMES:
     1: PN: US9958879 TABLE: 5 claimed protein
CN
CN
     3: PN: WOO0\$0587 SEQID: 1 claimed protein
CN
     Bio-Tropin
CN
     CB 311
CN
     Corpormon
CN
     Crescormon
CN
     Genotropin
     Growth hormone (human pituitary)
CN
```

```
CN
     Human growth hormone
CN
     Humatrope
     Infithopin CR
CN
     LY 137998
CN
     Nordit topin
CN
CN
     Norditropin SimpleXx
     Nordotrbpin
CN
CN
     Nutropin
CN
     Nutropin Depot
CN
     Saizen
CN
     SJ 0011
CN
     Somatogen\
     Somatotropin (human)
CN
     Somatropin'
CN
CN
     SR 29001
CN
     Valtropin
CN
     Zomacton
CN
     Zorbtive
     PROTEIN SEQUENCE
FS
     869741-23-1, 1145-52-1
DR
MF
     Unspecified
CI
     MAN
LC
     STN Files:
                  AMISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BIOSIS, BIOTECHNO,
       CA, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU,
       DRUGU, EMBASE \ IMSCOSEARCH, IMSDRUGNEWS, IMSPATENTS, IMSRESEARCH, IPA,
       MEDLINE, MRCK* PATDPASPC, PHAR, PIRA, PROMT, RTECS*, SCISEARCH,
       TOXCENTER, USAN, USPAT2, USPATFULL, VETU
         (*File contains numerically searchable property data)
     Other Sources:
                     LINECS**, WHO
         (**Enter CHEML\(TST File for up-to-date regulatory information)
**RELATED SEQUENCES AVAILABLE WITH SEQLINK**
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
*** USE 'SQD' OR 'SQIDE' FORMATS TO DISPLAY SEQUENCE ***
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
            1400 REFERENCES IN FILE CA. (1907 TO DATE)
              99 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
            1400_REFERENCES IN FILE CAPLUS (1907 TO DATE)
/L3
     ANSWER 3 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN
RN
     9050-36-6 REGISTRY
ÈΟ
     Entered STN: 16 Nov 1984
CN
     Maltodextrin (9CI) (CA INDEX NAME)
OTHER NAMES:
     Actistar 11700
CN
CN
     Amidex DE 10
CN
     C Pharm 01980
CN
     C*De Light 01970
CN
     C*deLight F 01970
CN
     C*deLight MD 01970
     C-Pur 01910
CN
     C-PUR 01915
CN
     C-Pur 01921
CN
CN
     C-Sperse MD 01314
     Cerestar C*PUR 01915
CN
CN
     Cerestar PUR 01915
CN
     DE 2
CN
     Dextrin, malto
CN
     Dry Sweet
CN
     Fibersol 2(E)
CN
     Foodtex
CN
     Frodex 10
CN
     Frodex 20
CN
     Glister
CN
     Glucidex 12
CN
     Glucidex 17
```

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CN
     Glucidex 19
CN
     Glucidex 19FD
CN
     Glucidex 2
     Glucidex 21
CN
     Glucidex 2B
CN
CN
     Glucidex 39
CN
     Glucidex 6
CN
     Glucidex IT 12
CN
     Glucidex IT 19
CN
     Glucidex IT 6
CN
     Glucidex IT 8
CN
     Instant N-Oil II
CN
     Instant Oil II
CN
     Instant Stellar
CN
     K 8
     Lodex 10
CN
CN
     Lodex 5
CN
     Lycadex 100
CN
     Lycadex 200
CN
     Lycatab
CN
     M 01960
CN
     M 040
CN
     Maldex 15
CN
     Maldex 150
CN
     Maldex 20
CN
     Maldex 30
CN
     Malta-Gran 10
CN
     Malta-Gran TG
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
     DISPLAY
DR
     126776-44-1, 126776-45-2, 127120-90-5, 54077-26-8, 104859-39-4,
     104859-43-0, 104859-45-2, 104859-47-4, 104859-49-6, 104859-62-3,
     104859-75-8, 61008-41-1, 142583-82-2, 89750-26-5, 87090-11-7, 39283-25-5,
     52769-80-9, 216252-89-0, 220857-34-1, 287179-53-7
MF
     Unspecified
CI
     PMS, COM, MAN
PCT
     Manual registration
LC
     STN Files:
                  AGRICOLA, ANABSTR, BIOSIS, BIOTECHNO, CA, CABA, CAPLUS,
       CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU,
       EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MSDS-OHS, PIRA, PROMT,
       TOXCENTER, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
                      DSL**, EINECS**, TSCA**
     Other Sources:
         (**Enter CHEMLIST File for up-to-date regulatory information)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
            3496 REFERENCES IN FILE CA (1907 TO DATE)
             154 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
            3507 REFERENCES IN FILE CAPLUS (1907 TO DATE)
     ANSWER 4 OF 10 REGISTRY
                                COPYRIGHT 2006 ACS on STN
\mathbf{L}_{\mathbf{Z}}
RN
     3040-38-8 REGISTRY
ΕD
     Entered STN: 16 Nov 1984
     1-Propanaminium, 2-(acetyloxy)-3-carboxy-N,N,N-trimethyl-, inner salt,
CN
     (2R) - (9CH) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN
     1-Propanaminium, 2-(acetyloxy)-3-carboxy-N,N,N-trimethyl-, inner salt,
CN
     Ammonium, (3-carboxy-2-hydroxypropyl)trimethyl-, hydroxide, inner salt,
     acetate, L- (8CI)
OTHER NAMES:
CN
     (-)-Acetylcarnitine
CN
     (R)-Acetylcarnitine
CN
     Acetyl-L-(-)-carnitine
CN
     Acetyl-L-carnitine
CN
     Acetylcarnitine
```

```
CN
     ALCAR
CN
     L-Acetylcarnitine
CN
     L-Carnitine acetyl ester
CN
     L-O-Acetylcarnitine
CN
     Levocarnitine acetyl
CN
     Nicetile
CN
     O-Acetyl-L-carnitine
     O-Acetylcarnitine
CN
FS
     STEREOSEARCH
DR
     461-77-8, 541-68-4, 3624-25-7, 74832-89-6
MF
CI
LC
     STN Files:
                   ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS,
       BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN,
       CSCHEM, DDFU, DRUGU, EMBASE, IMSCOSEARCH, IMSDRUGNEWS, IMSRESEARCH, IPA,
       MRCK*, PROMT, PROUSDDR, RTECS*, TOXCENTER, USPAT2, USPATFULL
          (*File contains numerically searchable property data)
     Other Sources:
Absolute stereochemistry.
                N+Me3
-02C
          OAc
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
             910 REFERENCES IN FILE CA (1907 TO DATE)
              20 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
              911 REFERENCES IN FILE CAPLUS (1907 TO DATE)
               3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
     ANSWER 5 OF 10 REGISTRY
                                COPYRIGHT 2006 ACS on STN
1.3
RN
     657-27-2 REGISTRY
     Entered STN: 16 Nov 1984
EE
     L-Lysine, monohydrochloride (9CI) (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
     Lysine, monohydrochloride, L- (8CI)
CN
OTHER NAMES:
CN
     Darvyl
CN
     L-Gen
CN
     L-Lysine hydrochloride
CN
     Lyamine
     Lysine hydrochloride
CN
     Lysine monohydrochloride
CN
CN
     Lysion
CN
     NSC 9253
FS
     STEREOSEARCH
     305-76-0, 93394-22-0
DR
     C6 H14 N2 O2 . Cl H
MF
CI
     COM
LC
     STN Files:
                   AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS,
       CASREACT, CBNB, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DETHERM*, EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MRCK*, MSDS-OHS,
       PATDPASPC, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN, USPAT7, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
CRN
     (56-87-1)
```

$$^{\text{NH}_2}$$
 $^{\text{NH}_2}$
 $^{\text{NH}_2}$
 $^{\text{NH}_2}$

HCl

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1118 REFERENCES IN FILE CA (1907 TO DATE)
31 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1118 REFERENCES IN FILE CAPLUS (1907 TO DATE)

ANSWER 6 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN

502-65-8 REGISTRY

ED Entered STN: 16 Nov 1984

CN ψ , ψ -Carotene (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Lycopene, all-trans- (8CI)

OTHER NAMES:

L3 RN

CN

CN (6E,8E,10E,12E,14E,16E,18E,20E,22E,24E,26E)-2,6,10,14,19,23,27,31-Octamethyl-2,6,8,10,12,14,16,18,20,22,24,26,30-dotriacontatridecaene

CN (all-E)-2,6,10,14,19,23,27,31-Octamethyl-2,6,8,10,12,14,16,18,20,22,24,26, 30-dotriacontatridecaene

2,6,8,10,12,14,16,18,20,22,24,26,30-Dotriacontatridecaene, 2,6,10,14,19,23,27,31-octamethyl-, (6E,8E,10E,12E,14E,16E,18E,20E,22E,24E, 26E)-

CN all-trans-Lycopene

CN C.I. 75125

CN Lyco Vit

CN Lycopene

CN Lycopene 7

CN NSC 407322

CN trans-Lycopene

FS STEREOSEARCH

DR 7634-65-3, 25453-98-9

MF C40 H56

CI COM

LC

STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IMSDRUGNEWS, IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, PIRA, PROMT, RTECS*, SCISEARCH, SPECINFO, TOXCENTER, USPAT2, USPATFULL (*File contains numerically searchable property data)

Other Sources: EINECS**

(**Enter CHEMLIST File for up-to-date regulatory information)

Double bond geometry as shown.

PAGE 1-A

PAGE 1-B

58: PN: US20040014159 SEQID: 33 unclaimed sequence

Benzenepropanoic acid, α -amino-4-hydroxy-, (S)-

CN CN

CN

CN

CN

L-(-)-Tyrosine

L-Phenylalanine, 4-hydroxy-

L-p-Tyrosine

```
CN
     NSC 82624
     NSC 9973
CN
CN
     p-Tyrosine
     Propanoic acid, 2-amino-3-(4-hydroxyphenyl)-, (S)-
CN
CN
     Tyrosine
FS
     STEREOSEARCH
     140-43-2, 55520-40-6, 1991-85-1, 46209-14-7
DR
MF
     C9 H11 N O3
CI
     COM
     STN Files:
                  ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS,
LC
       BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU,
       EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
       MSDS-OHS, NAPRALERT, PATDPASPC, PHAR, PIRA, PROMT, PS, RTECS*, SPECINFO,
       SYNTHLINE, TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, VETU
         (*File contains numerically searchable property data)
                     DSL**, EINECS**, TSCA**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)
Absolute stereochemistry. Rotation (-).
                   ∠CO2H
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
           43928 REFERENCES IN FILE CA (1907 TO DATE)
            1165 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
           43980 REFERENCES IN FILE CAPLUS (1907 TO DATE)
               7 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
L3
     ANSWER 9 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN
     56-85-9 REGISTRY
RN
     Entered STN: 16 Nov 1984
ED
     L-Glutamine (9CI) (CA INDEX NAME)
CN
OTHER CA INDEX NAMES:
CN
     Glutamine, L- (8CI)
OTHER NAMES:
CN
     (S)-2,5-Diamino-5-oxopentanoic acid
CN
     γ-Glutamine
CN
     2-Aminoglutaramic acid
CN
     Acustasin
CN
     Aesgen 14
CN
     Cebrogen
CN
     Glumin
     Glumin (amino acid)
CN
CN
     Glutamax
CN
     Glutamic acid 5-amide
     Glutamic acid amide
CN
CN
     Glutamine
CN
     L-(+)-Glutamine
CN
     L-2-Aminoglutaramidic acid
CN
     L-Glutamic acid \gamma-amide
CN
     Levoglutamide
CN
     NSC 27421
CN
     Pentanoic acid, 2,5-diamino-5-oxo-, (S)-
CN
     Stimulina
FS
     STEREOSEARCH
DR
     32640-56-5
MF
     C5 H10 N2 O3
CI
     COM
```

ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS,

BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS,

CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE,

LC

STN Files:

GMELIN*, IFICDB, IFIPAT, IFIUDB, IMSDRUGNEWS, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, PROMT, PS, RTECS*, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL

(*File contains numerically searchable property data)

(*File contains numerically searchable property data)

Other Sources: DSL**, EINECS**, TSCA**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

24245 REFERENCES IN FILE CA (1907 TO DATE)

474 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

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24276 REFERENCES IN FILE CAPLUS (1907 TO DATE)
                6 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
     ANSWER 10 OF 10 REGISTRY COPYRIGHT 2006 ACS on STN
L3
     50-23-7 REGISTRY
RN
ED
     Entered STN: 16 Nov 1984
     Pregn-4-ene-3,20-dione, 11,17,21-trihydroxy-, (11\beta)- (9CI) (CA INDEX
CN
     NAME)
OTHER CA INDEX NAMES:
     Cortisol (8CI)
CN
OTHER NAMES:
     11\beta, 17, 21-Trihydroxypregn-4-ene-3, 20-dione
CN
CN
     11β, 17, 21-Trihydroxyprogesterone
CN
     11\beta, 17\alpha, 21-Trihydroxypregn-4-ene-3, 20-dione
CN
     11β-Hydroxycortisone
CN
     17-Hydroxycorticosterone
CN
     17\alpha-Hydroxycorticosterone
CN
     4-Pregnene-11\beta, 17\alpha, 21-triol-3, 20-dione
CN
     Acticort
CN
     Aeroseb HC
     Ala-Cort
CN
CN
     Anflam
     Anti-inflammatory hormone
CN
CN
     CaldeCort Spray
CN
     CCN 90306A
CN
     Cetacort
CN
     Cobadex
```

CN Cortenema
CN Corticreme
CN Cortifan
CN Cortiment
CN Cortispray

Cort-Dome

Cortanal

Cortef

CN Cortonema
CN Cortril

CN

CN

CN

CN Cortril
CN Dermacort
CN Dermocortal
CN Dermolate

CN Dermolate CN Dihydrocostisone

CN Dinydrocostisone
CN Dioderm
CN Domolene-HC

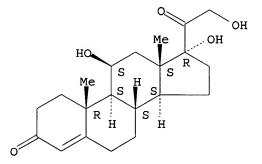
CN Domolene-HC CN Efcorbin CN Efcortelan

CN Eldecort CN Epiderm H

```
CN
     Ficortril
CN
     Genacort
CN
     HC
CN
     Heb-Cort
CN
    Hidro-Colisona
CN
    Hycort
CN
    Hycortol
CN
    Hycortole
CN
     Hydracort
CN
     Hydrasson
     Hvdro-Adreson
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
     DISPLAY
FS
     STEREOSEARCH
DR
     8056-08-4, 8063-42-1, 80562-38-5
MF
     C21 H30 O5
     COM
CI
     STN Files:
                 ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS,
LC
      BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU,
      EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSDRUGNEWS, IMSPATENTS,
       IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, PHAR, PIRA,
       PROMT, PS, RTECS*, SCISEARCH, SPECINFO, SYNTHLINE, TOXCENTER, USAN,
      USPAT2, USPATFULL, VETU
         (*File contains numerically searchable property data)
    Other Sources:
                     DSL**, EINECS**, TSCA**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)
```

CN

Esiderm H Evacort



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

37291 REFERENCES IN FILE CA (1907 TO DATE)

351 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

37322 REFERENCES IN FILE CAPLUS (1907 TO DATE)

20 REFERENCES IN FILE CAOLD (PRIOR TO 1967)